Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112221\

Data File : BN017544.D

Acq On : 22 Nov 2021 13:26

Operator : CG/JU Sample : SSTD16050

Misc :

ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 22 16:05:58 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

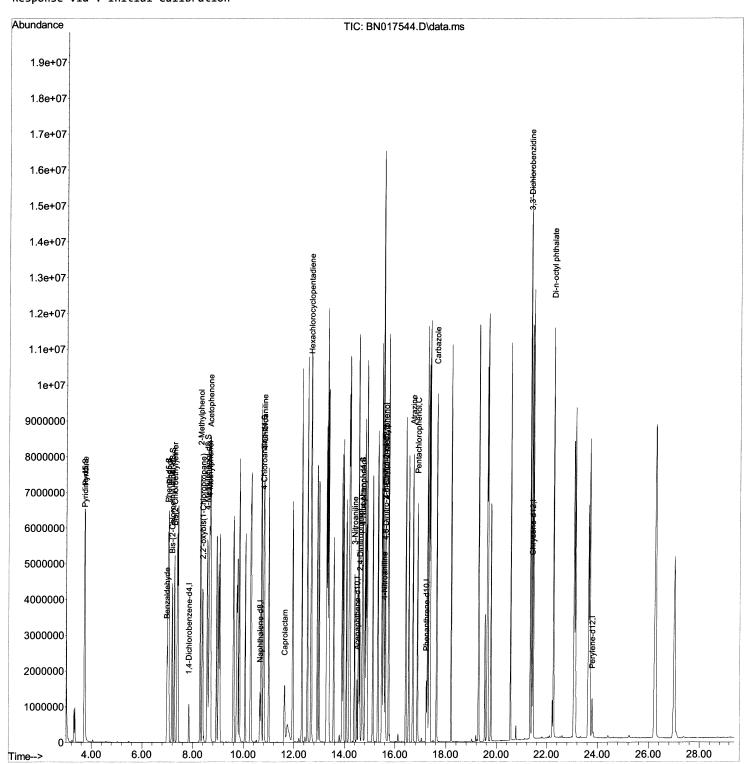
Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 15:40:34 2021

Response via : Initial Calibration

Instrument : BNA_N ClientSampleId : SSTD160250

Manual IntegrationsAPPROVED



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112221\

Data File : BN017544.D Acq On : 22 Nov 2021 13:26

Operator : CG/JU Sample : SSTD16050

Misc :

ALS Vial : 7 Sample Multiplier: 1

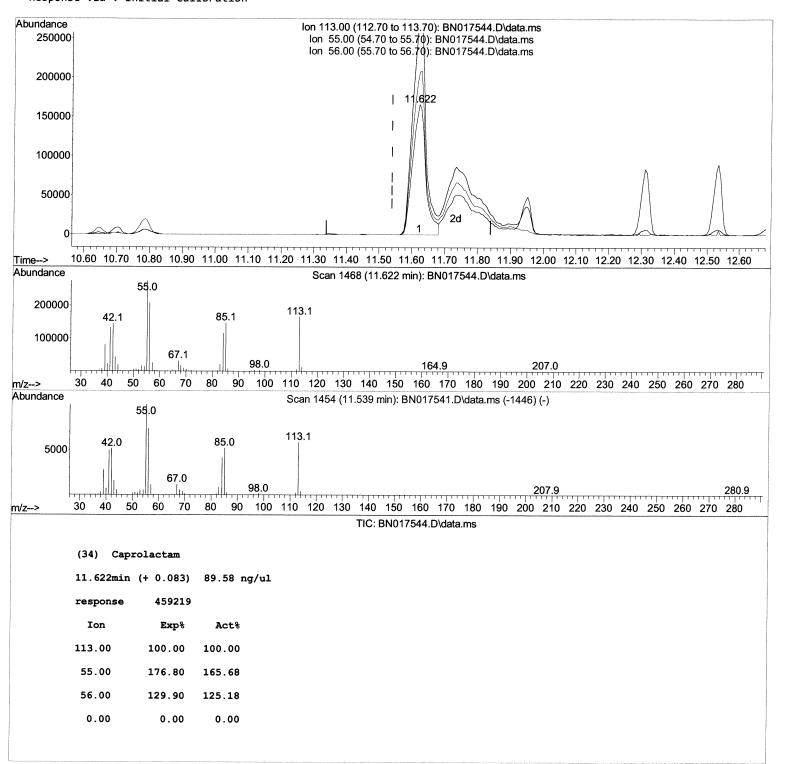
Quant Time: Nov 22 16:05:58 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN112221.M

Quant Title : SVOA CALIBRATION

QLast Update : Mon Nov 22 15:40:34 2021 Response via : Initial Calibration Instrument : BNA_N ClientSampleld : SSTD160250

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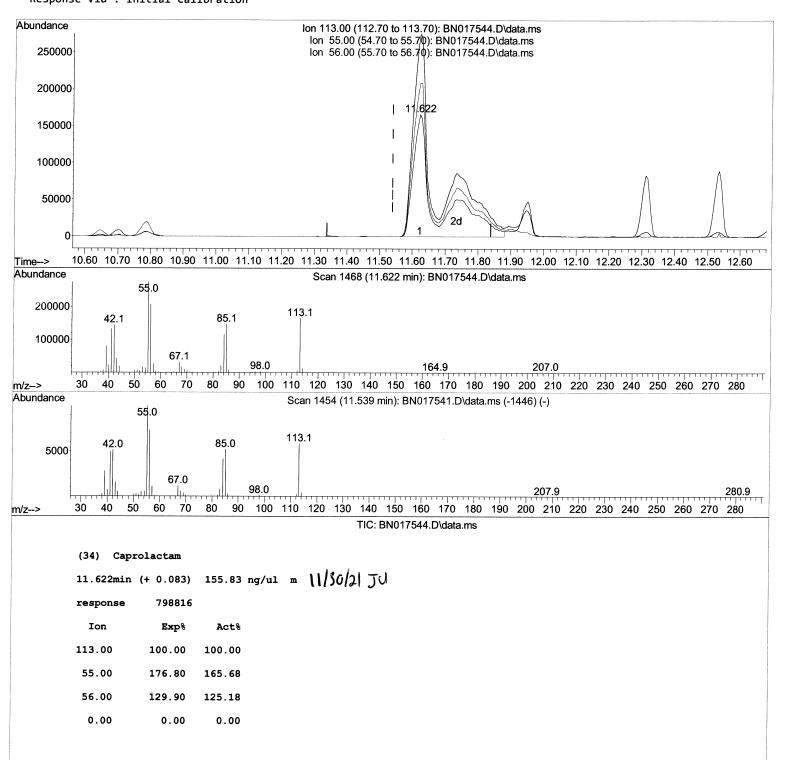
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 $\label{lem:quant_Method} {\tt Quant_Methods\SFAM-EPA-BN112221.M}$

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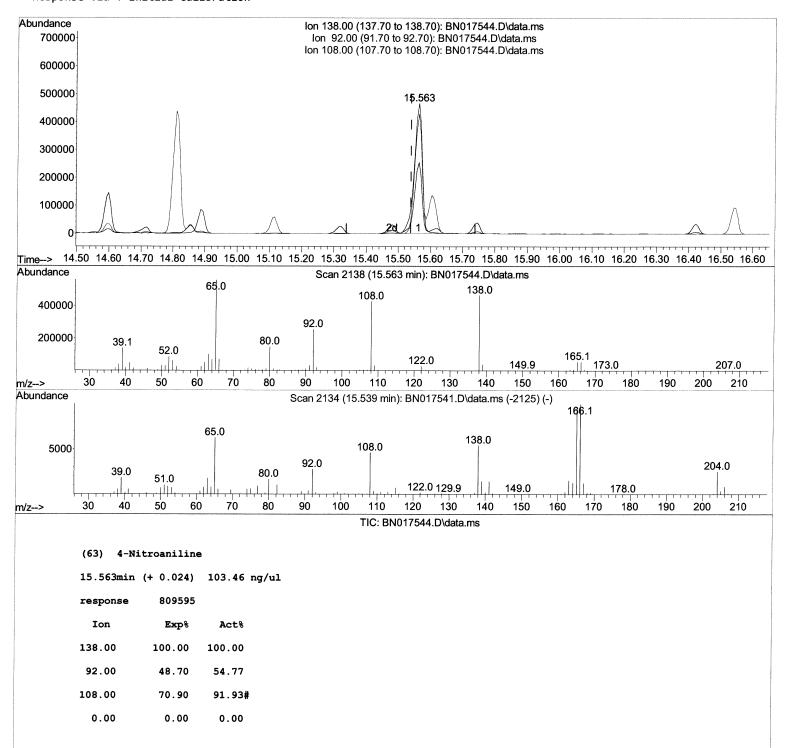
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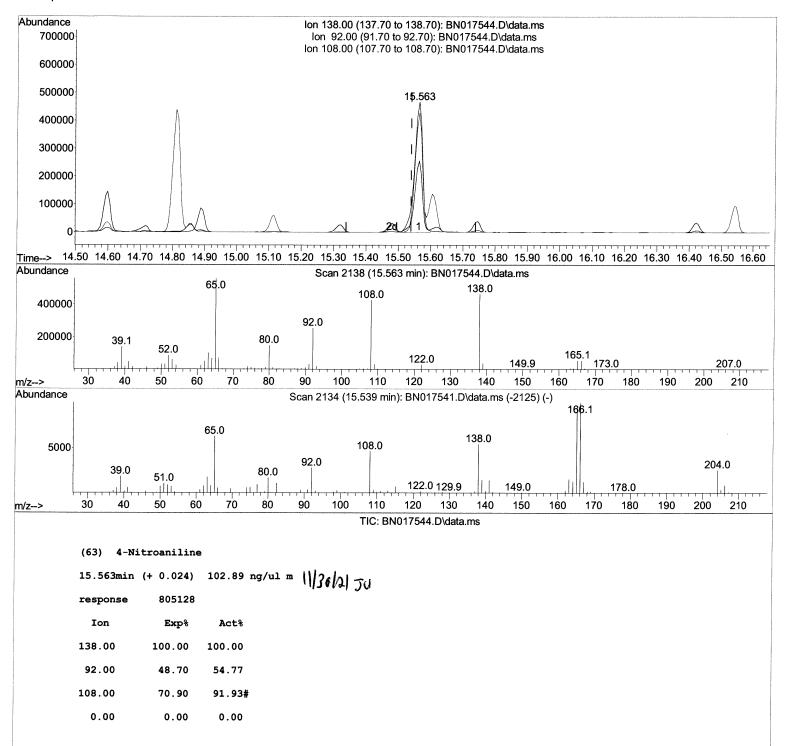
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Manual IntegrationsAPPROVED

Compound	R.T.	QIon	Response	Conc Un	its Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	270887	20 000	ng/ul	0.00
20) Naphthalene-d8	10.646		1133645		ng/ul	0.00
38) Acenaphthene-d10	14.487		622972		ng/ul	0.00
64) Phenanthrene-d10	17.228		1048699		ng/ul	0.00
79) Chrysene-d12	21.416		675407		ng/ul	0.00
88) Perylene-d12	23.786		719513		ng/ul	0.01
oo) Terytene utz	23.700	204	,15515	20.000	iig/ui	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.711	84	3011552	151.413	<u> </u>	0.00
7) Phenol-d5	7.028	99	3754456	152.245	_	0.02
<pre>9) Bis-(2-Chloroethyl)eth</pre>	7.187	67	2221048	143.724	-	0.01
11) 2-Chlorophenol-d4	0.000	132	0d		ng/ul	
15) 4-Methylphenol-d8	8.569	113	2829508	145.868		0.02
21) Nitrobenzene-d5	0.000	128	0d		ng/ul	
24) 2-Nitrophenol-d4	0.000	143	Ød		ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d		ng/ul	
31) 4-Chloroaniline-d4	10.787	131	3267397	136.018	_	0.02
46) Dimethylphthalate-d6	0.000	166	0d		ng/ul	
49) Acenaphthylene-d8	0.000	160	9d		ng/ul	
54) 4-Nitrophenol-d4	14.698	143	1130958	147.553	_	0.04
60) Fluorene-d10	0.000	176	0d		ng/ul	
65) 4,6-Dinitro-2-methylph	15.604	200	924839	169.793	-	0.02
73) Anthracene-d10	0.000	188	0d		ng/ul	*****
81) Pyrene-d10	0.000	212	0d		ng/ul	
92) Benzo(a)pyrene-d12	0.000	264	0d		ng/ul	
Target Compounds					Qva.	luo
5) Pyridine	3.728	79	2970823	148.424	-	
6) Benzaldehyde	6.981	73 77	943206	72.101	_	97 92
8) Phenol	7.058	94	3665748	146.769		93
10) Bis(2-Chloroethyl)ether	7.038	93	2896930	141.458		98 98
13) 2-Methylphenol	8.299					99
14) 2,2'-oxybis(1-Chloropr	8.387	108 45	2741687 3867326	145.744 140.072		99 96
16) Acetophenone	8.681	105	3866297		_	
18) 4-Methylphenol	8.652	108	2862679	133.282 142.053		94 97
32) 4-Chloroaniline	10.810	127	3186225		-	99
34) Caprolactam	11.622	113	798816m≯	130.726		
40) Hexachlorocyclopentadiene	12.669	237	1897869	160.144		11/36/2174
51) 3-Nitroaniline					-	
53) 2,4-Dinitrophenol	14.393 14.598	138	1166566 780705	130.774		88
55) 4-Nitrophenol		184		184.319		93 76
63) 4-Nitrophenol	14.716	109	988080	147.805	_	76
66) 4,6-Dinitro-2-methylph	15.563	138	805128m>		_	1112012124
70) Atrazine	15.622	198	892109	163.136		83
71) Pentachlorophenol	16.704	200	1622544	145.338	-	98
71) Pentachiorophenoi 77) Carbazole	16.881	266 167	1035380	165.011		100
84) 3,3'-Dichlorobenzidine	17.634	167	6534117	134.173	•	99
89) Di-n-octyl phthalate	21.333 22.257	252 149	1737968 8261416	135.270 151.178		93 100
					-	

Quantitation Report (QT Reviewed)

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Compound R.T. QIon Response Conc Units Dev(Min)

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Instrument:
BNA_N
ClientSampleId:
SSTD160250

Manual IntegrationsAPPROVED